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PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

- NEWS 1 Web Page for STN Seminar Schedule N. America NEWS 2 DEC 01 ChemPort single article sales feature unavailable NEWS 3 APR 03 CAS coverage of exemplified prophetic substances enhanced
- NEWS 4 APR 07 STN is raising the limits on saved answers
- NEWS 5 APR 24 CA/Caplus now has more comprehensive patent assignee information
- NEWS 6 APR 26 USPATFULL and USPAT2 enhanced with patent assignment/reassignment information
- NEWS 7 APR 28 CAS patent authority coverage expanded
- NEWS 8 APR 28 ENCOMPLIT/ENCOMPLIT2 search fields enhanced
- NEWS 9 APR 28 Limits doubled for structure searching in CAS REGISTRY
- NEWS 10 MAY 08 STN Express, Version 8.4, now available
- NEWS 11 MAY 11 STN on the Web enhanced
- NEWS 12 MAY 11 BEILSTEIN substance information now available on STN Easy
- NEWS 13 MAY 14 DGENE, PCTGEN and USGENE enhanced with increased limits for exact sequence match searches and introduction of free HIT display format
- NEWS 14 MAY 15 INPADOCDB and INPAFAMDB enhanced with Chinese legal status data
- NEWS 15 MAY 28 CAS databases on STN enhanced with NANO super role in records back to 1992
- NEWS 16 JUN 01 CAS REGISTRY Source of Registration (SR) searching enhanced on STN
- NEWS 17 JUN 26 NUTRACEUT and PHARMAML no longer updated
- NEWS 18 JUN 29 IMSCOPROFILE now reloaded monthly
- NEWS 19 JUN 29 EPFULL adds SLART to AB, MCLM, and TI fields
- NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4, AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.
- NEWS HOURS STN Operating Hours Plus Help Desk Availability NEWS LOGIN Welcome Banner and News Items

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FILE 'HOME' ENTERED AT 13:08:14 ON 01 JUL 2009

=> file reg

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
0.22
0.22

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STRUCTURE FILE UPDATES: 30 JUN 2009 HIGHEST RN 1160555-05-4 DICTIONARY FILE UPDATES: 30 JUN 2009 HIGHEST RN 1160555-05-4

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http://www.cas.org/support/stngen/stndoc/properties.html

```
=> e 1-hydroxy-2,3-dihydroindole/cn
                   1 1-HYDROXY-2,3-DICARBOMETHOXY-1,3-CYCLOHEPTADIENE/CN
E1
E2
                    1
                            1-HYDROXY-2,3-DIETHOXYACRIDONE/CN
Е3
                    0 --> 1-HYDROXY-2,3-DIHYDROINDOLE/CN
                   1 1-HYDROXY-2,3-DIIODO-2-PROPENE/CN
1 1-HYDROXY-2,3-DIMETHOXY-10-METHYL
E4
E5
                           1-HYDROXY-2,3-DIMETHOXY-10-METHYLACRIDAN-9-ONE/CN
                  1 1-HYDROXY-2,3-DIMETHOXY-10-METHYLACRIDONE/CN
1 1-HYDROXY-2,3-DIMETHOXY-6-FORMYL-6A,7-DEHYDROAPORPHINE/CN
1 1-HYDROXY-2,3-DIMETHOXY-7-METHYL-9,10-ANTHRAQUINONE/CN
1 1-HYDROXY-2,3-DIMETHOXYACRIDAN-9-ONE/CN
1 1-HYDROXY-2,3-DIMETHOXYBENZENE/CN
1 1-HYDROXY-2,3-DIMETHOXYXANTHONE/CN
E.7
E9
E10
E11
                            1-HYDROXY-2,3-DIMETHYL-N-METHYLACRIDONE/CN
E12
```

=> logoff hold

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
0.96
1.18

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 13:09:40 ON 01 JUL 2009

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID: SSSPTA1623PA7

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * * SESSION RESUMED IN FILE 'REGISTRY' AT 13:14:08 ON 01 JUL 2009 FILE 'REGISTRY' ENTERED AT 13:14:08 ON 01 JUL 2009 COPYRIGHT (C) 2009 American Chemical Society (ACS)

COST IN U.S. DOLLARS

SINCE FILE TOTAL
ENTRY SESSION
0.96 1.18

11 ANSWERS

FULL ESTIMATED COST

=>

 $\begin{tabular}{ll} Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary files\10516979\10516979 dihydroindole.str \end{tabular}$

L1 STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STR

Structure attributes must be viewed using STN Express query preparation.

=> search la sss sam

SAMPLE IS IGNORED AS A SCOPE FOR THIS SEARCH

83493 LA

523 LAS

L2 83998 LA

(LA OR LAS)

=> search l1 sss sam

SAMPLE SEARCH INITIATED 13:17:32 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 4840 TO ITERATE

41.3% PROCESSED 2000 ITERATIONS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 92628 TO 100972 PROJECTED ANSWERS: 223 TO 841

L3 11 SEA SSS SAM L1

=> d scan

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 1H-Indole, 2,3-dihydro-1-hydroxy-2-[2-[3-[2-(1-hydroxy-1H-indol-2-yl)ethenyl]-2-cyclohexen-1-ylidene]ethylidene]-, perchlorate (1:1)

MF C26 H24 N2 O2 . C1 H O4

CM 1

CM 2

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):5

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN Hydrazinecarbothioamide, 2-(1,2-dihydro-1-hydroxy-2-oxo-3H-indol-3-ylidene)-

MF C9 H8 N4 O2 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 1H-Indol-3-ol, 4,5,6,7-tetrafluoro-2,3-dihydro-1-hydroxy-

MF C8 H5 F4 N O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 2-Thiophenecarboxamide, N-[2,3-dihydro-1-hydroxy-3-(1-hydroxypropylidene)-2-oxo-1H-indol-6-yl]-

MF C16 H14 N2 O4 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 1H-Indole-2,3-dione, 6-chloro-1-hydroxy-

MF C8 H4 Cl N O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 11 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

 $\label{eq:local_section} \mbox{IN} \quad \mbox{Benzenamine, N-(2-ethyl-1,2-dihydro-1-hydroxy-2-phenyl-3H-indol-3-ylidene)-1} \\ \mbox{Benzenamine, N-(2-ethyl-1,2-dihydro-1-hydroxy-2-phenyl-3-hydroxy-2-phenyl-3-hydroxy-2-p$

MF C22 H20 N2 O

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> search 11 sss full FULL SEARCH INITIATED 13:18:09 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 98117 TO ITERATE

100.0% PROCESSED 98117 ITERATIONS SEARCH TIME: 00.00.02

304 ANSWERS

L4 304 SEA SSS FUL L1

=> save temp 14 polyinhib/a ANSWER SET L4 HAS BEEN SAVED AS 'POLYINHIB/A'

accessing the remaining file names entered.

=> file caplus\
'CAPLUS\' IS NOT A VALID FILE NAME
SESSION CONTINUES IN FILE 'REGISTRY'
Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue

=> file caplus
COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 195.55 195.77

FILE 'CAPLUS' ENTERED AT 13:19:02 ON 01 JUL 2009
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FILE COVERS 1907 - 1 Jul 2009 VOL 151 ISS 1 FILE LAST UPDATED: 30 Jun 2009 (20090630/ED) REVISED CLASS FIELDS (/NCL) LAST RELOADED: Apr 2009 USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Apr 2009

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 15

L5 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> 14

L5 244 L4

=> polymeriz?

423002 POLYMERIZ?

96532 POLYMD

96532 POLYMD

37642 POLYMG

386985 POLYMN

10455 POLYMNS

388377 POLYMN

(POLYMN OR POLYMNS)

L6 611277 POLYMERIZ?

(POLYMERIZ? OR POLYMD OR POLYMG OR POLYMN)

=> 15 and 16

L7 2 L5 AND L6

=> d 17 1-2 ti fbib abs

L7 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN

TI Abilities of 1-(9-anthrylmethyloxy)-2-pyridone and related compounds to initiate radical and cationic photopolymerizations

AN 2004:474370 CAPLUS <<LOGINID::20090701>>

DN 141:157530

TI Abilities of 1-(9-anthrylmethyloxy)-2-pyridone and related compounds to initiate radical and cationic photopolymerizations

AU Tanaka, Kenta; Nakamura, Kanako; Yoshioka, Nariyoshi; Kameyama, Atsushi; Igarashi, Tetsutaro; Sakurai, Tadamitsu

CS High-Tech Research Center, Kanagawa University, Yokohama, 221-8686, Japan

SO Journal of Polymer Science, Part A: Polymer Chemistry (2004), 42(12), 2859-2865

CODEN: JPACEC; ISSN: 0887-624X

PB John Wiley & Sons, Inc.

DT Journal

LA English

AB This study explored the abilities of 1-(9-anthrylmethyloxy)-2-pyridone and related compds., which absorb long-wavelength light (>350 nm), to photochem. initiate radical and cationic polymns. It was found that the irradiation of the title compds. initiates the radical polymn. of styrene whereas the cationic polymerization of oxetane proceeds in the presence of these photoinitiators to a negligible extent. The

behavior of 9-anthrylmethyloxyl and amidyl radicals in the photopolymn. process of styrene was discussed based on 1H NMR, UV, and fluorescence spectral data. In addition, the photoinitiation ability of the anthrylmethyloxyl end group was also investigated by using its model compound

RE.CNT 16 THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

- L7 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN
- TI Use of cyclic hydroxylamines as polymerization inhibitors
- AN 2003:1006927 CAPLUS <<LOGINID::20090701>>
- DN 140:43124
- TI Use of cyclic hydroxylamines as polymerization inhibitors
- IN Philips, Emyr; Loyns, Colin
- PA A H Marks & Company Limited, UK
- SO PCT Int. Appl., 19 pp. CODEN: PIXXD2
- DT Patent
- LA English
- FAN.CNT 1

r AIN.					KIND DATE		APPLICATION NO.					DATE						
ΡI	WO	2003	03106390 A1 20031224 WO 2003-GB2367															
		W:	CO, GM, LS,	CR, HR, LT,	CU, HU, LU,	CZ, ID, LV,	DE, IL, MA,	DK, IN, MD,	DM, IS, MG,	DZ, JP, MK,	EC, KE, MN,	BG, EE, KG, MW,	ES, KP, MX,	FI, KR, MZ,	GB, KZ, NO,	GD, LC, NZ,	GE, LK, OM,	GH, LR, PH,
		RW:	UA,	UG,	US,	UZ,	VC,	VN,	YU,	ZA,	ZM,	SL, ZW TZ,						
			KG, FI,	KZ, FR,	MD, GB,	RU, GR,	TJ, HU,	TM, IE,	AT, IT,	BE, LU,	BG, MC,	CH, NL, GW,	CY, PT,	CZ, RO,	DE, SE,	DK, SI,	EE, SK,	ES, TR,
											GB 2 GB 2	2002- 2002-	1348 1409	0		A 2 A 2	0020 0020	613 619
	AU	2003	2447	50		A1		2003	1231		GB 2 GB 2	2003- 2002- 2002- 2003-	1348 1409	0 3		A 2 A 2	0020	613 619
	EP	1511	704			A1		2005	0309		EP 2	2003-	7382	35		2	0030	530
		R:		SI,	LT,	LV,	FI,			CY,	AL, GB 2 GB 2	IT, TR, 2002- 2002- 2003-	BG, 1348 1409	CZ, 0 3	EE,	HU, A 2 A 2	SK 0020	613 619
	JP	2005529223			Τ	20050929		JP 2004-513226 GB 2002-13480 GB 2002-14093 WO 2003-GB2367				A 2 A 2	0030 0020 0020 0030	613 619				
	US	2006	0167	244		A1		2006	0727		GB 2 GB 2	2005- 2002- 2002- 2003-	1348 1409	0 3		A 2 A 2	0050 0020 0020 0030	613 619

OS MARPAT 140:43124

AB A polymerization inhibitor comprising a non-hindered cyclic hydroxylamine (e.g. 1-hydroxypiperidine) either alone or in combination with an addnl. inhibitor is described for a variety of monomers (e.g., styrene).

RE.CNT 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 12.74 208.51

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE TOTAL
ENTRY SESSION
-1.64 -1.64

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http://www.cas.org/support/stngen/stndoc/properties.html

=>

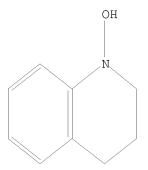
 $\begin{tabular}{l} Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary files\10516979\trihydroquinoline.str \end{tabular}$

L8 STRUCTURE UPLOADED

=> d 18

L8 HAS NO ANSWERS

L8 STR



Structure attributes must be viewed using STN Express query preparation.

=> search 18 sss sam
SAMPLE SEARCH INITIATED 13:25:03 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 4840 TO ITERATE

41.3% PROCESSED 2000 ITERATIONS 5 ANSWERS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 92628 TO 100972
PROJECTED ANSWERS: 34 TO 450

L9 5 SEA SSS SAM L8

=> d scan

L9 5 ANSWERS REGISTRY COPYRIGHT 2009 ACS on STN

IN 2(1H)-Quinolinone, 3-amino-3,4-dihydro-1-hydroxy-, hydrochloride (1:1)

MF C9 H10 N2 O2 . C1 H

● HCl

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> search 18 sss full FULL SEARCH INITIATED 13:25:36 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 98117 TO ITERATE

100.0% PROCESSED 98117 ITERATIONS 333 ANSWERS

SEARCH TIME: 00.00.02

L10 333 SEA SSS FUL L8

=> save temp 110 quins/a
ANSWER SET L10 HAS BEEN SAVED AS 'QUINS/A'

=> file caplus

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 186.84 395.35

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL

ENTRY SESSION

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FILE COVERS 1907 - 1 Jul 2009 VOL 151 ISS 1
FILE LAST UPDATED: 30 Jun 2009 (20090630/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Apr 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Apr 2009

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=> 110

L11 215 L10

=> dhis

L12 30 DHIS

=> d his

T.1

(FILE 'HOME' ENTERED AT 13:08:14 ON 01 JUL 2009)

FILE 'REGISTRY' ENTERED AT 13:08:24 ON 01 JUL 2009 E 1-HYDROXY-2,3-DIHYDROINDOLE/CN

STRUCTURE UPLOADED

L2 83998 SEARCH LA SSS SAM

L3 83998 SEARCH LA SSS SAM L3 11 SEARCH L1 SSS SAM

L4 304 SEARCH L1 SSS FULL

SAVE TEMP L4 POLYINHIB/A

FILE 'CAPLUS' ENTERED AT 13:19:02 ON 01 JUL 2009

L5 244 L4

L6 611277 POLYMERIZ?

L7 2 L5 AND L6

FILE 'REGISTRY' ENTERED AT 13:24:19 ON 01 JUL 2009

L8 STRUCTURE UPLOADED

L9 5 SEARCH L8 SSS SAM

L10 333 SEARCH L8 SSS FULL

SAVE TEMP L10 QUINS/A

FILE 'CAPLUS' ENTERED AT 13:26:08 ON 01 JUL 2009

L11 215 L10 L12 30 DHIS

=> file 16 and 111

'L6' IS NOT A VALID FILE NAME

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ENTER A FILE NAME OR (IGNORE):end

=> 16 and 111

L13 4 L6 AND L11

=> d 113 1-4 ti fbib abs

- L13 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN
- TI Pharmacophore based discovery of potential antimalarial agent targeting haem detoxification pathway
- AN 2008:1462479 CAPLUS <<LOGINID::20090701>>
- DN 150:113707
- TI Pharmacophore based discovery of potential antimalarial agent targeting haem detoxification pathway
- AU Acharya, Badri Narayan; Saraswat, Deepika; Kaushik, Mahabir Parshad
- CS Discovery Centre, Defence Research and Development Establishment, Madhya Pradesh, 474002, India
- SO European Journal of Medicinal Chemistry (2008), 43(12), 2840-2852 CODEN: EJMCA5; ISSN: 0223-5234
- PB Elsevier Masson SAS
- DT Journal
- LA English
- AB Pharmacophore hypotheses were generated from mols. having putative antimalarial activities targeting haem detoxification pathway of malarial parasite. A training set consisting of 33 compds., whose activities were evaluated for haem polymerization inhibition and against chloroquine resistant (K1) strain of Plasmodium falciparum, was optimized to generate hypotheses. The hypothesis showing optimum correlation between actual and estimated activities was validated by Fischer's randomization test at 95% confidence level and used as a model to screen our inhouse compound database. Nicotinic acid [trans-3-(4-ethoxy-3-methoxy-phenyl)-1-(4-hydroxy-phenyl)-allylidene]-hydrazide (ALH5) was obtained as a hit. The compound was synthesized and evaluated against chloroquine sensitive (MRC-02) and resistant (RKL9) strains of malarial parasite P. falciparum. The compound showed antimalarial activity in nanomolar range and was found comparable with chloroquine.
- RE.CNT 31 THERE ARE 31 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT
- L13 ANSWER 2 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN
- TI Pharmacophore-based predictive model generation for potent antimalarials targeting haem detoxification pathway
- AN 2008:322731 CAPLUS <<LOGINID::20090701>>
- DN 148:486299
- TI Pharmacophore-based predictive model generation for potent antimalarials targeting haem detoxification pathway
- AU Acharya, Badri Narayan; Kaushik, Mahabir Parshad
- CS Defence Research and Development Establishment, Discovery Center, Gwalior, 474002, India
- SO Medicinal Chemistry Research (2008), Volume Date 2007, 16(5), 213-229

CODEN: MCREEB; ISSN: 1054-2523

- PB Birkhaeuser Boston
- DT Journal
- LA English

AB Pharmacophore hypotheses were developed for mols. having antimalarial activities targeting the haem detoxification pathway of the malaria parasite. A training set consisting of 23 compds. was selected to generate these hypotheses, and their activities were evaluated for haem polymerization inhibition and against chloroquine-sensitive (3D7) as well as chloroquine-resistant (K1) strains of P. falciparum. The models were cross-validated by Fischer's randomization test at a 95% confidence level. The model developed against chloroquine-resistant malaria parasites was found to yield the best predictions among the three models.

RE.CNT 23 THERE ARE 23 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

- L13 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN
- TI Photosensitive composition sensitive to 350-450 nm laser for manufacturing CTP (Computer-to-Plate) lithographic printing plate
- AN 2004:931941 CAPLUS <<LOGINID::20090701>>
- DN 141:403505
- TI Photosensitive composition sensitive to 350-450 nm laser for manufacturing CTP (Computer-to-Plate) lithographic printing plate
- IN Ishiji, Yohei; Shibuya, Akinori
- PA Fuji Photo Film Co., Ltd., Japan
- SO Jpn. Kokai Tokkyo Koho, 54 pp. CODEN: JKXXAF
- DT Patent
- LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
ΡI	JP 2004309977	A	20041104	JP 2003-106678	20030410
				JP 2003-106678	20030410

OS MARPAT 141:403505

GΙ

- AB The title photosensitive composition comprises a photosensitizing dye represented by I (R1-9 = H, monovalent nonmetal group; Z = nonmetal atoms for completing ring), an activation compound capable of generating radical or acid, and a polymerizable compound capable of being polymerized upon contact with radical or acid.
- L13 ANSWER 4 OF 4 CAPLUS COPYRIGHT 2009 ACS on STN
- TI Use of cyclic hydroxylamines as polymerization inhibitors
- AN 2003:1006927 CAPLUS <<LOGINID::20090701>>

```
140:43124
DN
ТΤ
     Use of cyclic hydroxylamines as polymerization inhibitors
ΙN
     Philips, Emyr; Loyns, Colin
PA
     A H Marks & Company Limited, UK
SO
     PCT Int. Appl., 19 pp.
     CODEN: PIXXD2
DT
     Patent
     English
FAN.CNT 1
                        KIND DATE
                                           APPLICATION NO.
     PATENT NO.
                        ____
                                           _____
     WO 2003106390
                         A1 20031224 WO 2003-GB2367
                                                                   20030530
         W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
             CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
             GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH,
             PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ,
             UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,
             KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR,
             BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
                                             GB 2002-13480 A 20020613
GB 2002-14093 A 20020619
     AU 2003244750
                          Α1
                                20031231
                                             AU 2003-244750
                                                                     20030530
                                                                A 20020613
A 20020619
                                             GB 2002-13480
                                             GB 2002-14093
                                                                W 20030530
                                             WO 2003-GB2367
     EP 1511704
                          A1
                                20050309
                                             EP 2003-738235
                                                                     20030530
         R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
             IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK
                                             GB 2002-13480 A 20020613
                                             GB 2002-14093
                                                                A 20020619
                                             WO 2003-GB2367
                                                                W 20030530
     JP 2005529223
                          Τ
                                 20050929
                                             JP 2004-513226
                                                                     20030530
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OS MARPAT 140:43124

US 20060167244

AB A polymerization inhibitor comprising a non-hindered cyclic hydroxylamine (e.g. 1-hydroxypiperidine) either alone or in combination with an addnl. inhibitor is described for a variety of monomers (e.g., styrene).

20060727

GB 2002-13480

WO 2003-GB2367

WO 2003-GB2367 W 20030530

US 2005-516979

GB 2002-13480

GB 2002-14093

GB 2002-14093

A 20020613

A 20020619

W 20030530

A 20020613

A 20020619

20050810

RE.CNT 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

Α1

=> logoff hold COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY 17.74	SESSION 413.09
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-3.28	-4.92

STN INTERNATIONAL SESSION SUSPENDED AT 13:30:10 ON 01 JUL 2009

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 13:30:32 ON 01 JUL 2009 FILE 'CAPLUS' ENTERED AT 13:30:32 ON 01 JUL 2009 COPYRIGHT (C) 2009 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	17.74	413.09
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-3.28	-4.92

=>

=> save temp all ammendedsrch/l L# LIST L1-L13 HAS BEEN SAVED AS 'AMMENDEDSRCH/L'

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	33.24	428.59
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-3.28	-4.92

SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 13:49:24 ON 01 JUL 2009